

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061623\
 Data File : BN025884.D
 Acq On : 16 Jun 2023 15:35
 Operator : MA/JU
 Sample : 03225-02DL 10X
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 GWOW-BR-05-75-87-061323DL

Quant Time: Jun 16 16:45:41 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN053123.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 16 14:10:48 2023
 Response via : Initial Calibration

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 06/19/2023
 Supervised By :mohammad ahmed 06/20/2023

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.005	152	4693	0.400	ng	0.00
7) Naphthalene-d8	10.824	136	11734	0.400	ng	0.01
13) Acenaphthene-d10	14.630	164	6805	0.400	ng	0.00
19) Phenanthrene-d10	17.381	188	14377	0.400	ng	0.00
29) Chrysene-d12	21.560	240	10061	0.400	ng	# 0.00
35) Perylene-d12	24.045	264	10391	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.571	112	147	0.011	ng	0.03
5) Phenol-d6	7.196	99	100	0.006	ng	0.04
8) Nitrobenzene-d5	9.223	82	258	0.024	ng	0.05
11) 2-Methylnaphthalene-d10	12.434	152	561	0.033	ng	0.04
14) 2,4,6-Tribromophenol	16.152	330	72	0.034	ng	0.02
15) 2-Fluorobiphenyl	13.283	172	873	0.030	ng	0.02
27) Fluoranthene-d10	19.416	212	1393	0.047	ng	0.00
31) Terphenyl-d14	20.001	244	1063	0.044	ng	0.00
Target Compounds						
2) 1,4-Dioxane	3.383	88	19034	3.457	ng	# 93
34) Bis(2-ethylhexyl)phtha...	21.453	149	1173	0.051	ng	# 99
36) Indeno(1,2,3-cd)pyrene	26.703	276	910m	0.020	ng	
37) Benzo(b)fluoranthene	23.294	252	1382	0.034	ng	# 1
38) Benzo(k)fluoranthene	23.347	252	956	0.023	ng	# 1
41) Benzo(g,h,i)perylene	27.463	276	1009m	0.025	ng	

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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